

STN Columbus

COST IN U.S. DOLLARS	SINCE FILE	TOTAL
	ENTRY	SESSION
FULL ESTIMATED COST	0.21	0.21

FILE 'REGISTRY' ENTERED AT 14:01:36 ON 23 DEC 2003
 USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT.
 PLEASE SEE "HELP USAGETERMS" FOR DETAILS.
 COPYRIGHT (C) 2003 American Chemical Society (ACS)

Property values tagged with IC are from the ZIC/VINITI data file
 provided by InfoChem.

STRUCTURE FILE UPDATES: 22 DEC 2003 HIGHEST RN 630045-65-7
 DICTIONARY FILE UPDATES: 22 DEC 2003 HIGHEST RN 630045-65-7

TSCA INFORMATION NOW CURRENT THROUGH JULY 14, 2003

Please note that search-term pricing does apply when
 conducting SmartSELECT searches.

Crossover limits have been increased. See HELP CROSSOVER for details.

Experimental and calculated property data are now available. For more
 information enter HELP PROP at an arrow prompt in the file or refer
 to the file summary sheet on the web at:
<http://www.cas.org/ONLINE/DBSS/registryss.html>

=> s fluprostenol

L1 4 FLUPROSTENOL

=> d l1 4

L1 ANSWER 4 OF 4 REGISTRY COPYRIGHT 2003 ACS on STN

RN 40666-16-8 REGISTRY

CN 5-Heptenoic acid, 7-[(1R,2R,3R,5S)-3,5-dihydroxy-2-[(1E,3R)-3-hydroxy-4-[3-(trifluoromethyl)phenoxy]-1-butenyl]cyclopentyl]-, (5Z)-rel- (9CI) (CA INDEX NAME)

OTHER CA INDEX NAMES:

CN 5-Heptenoic acid, 7-[3,5-dihydroxy-2-[3-hydroxy-4-[3-(trifluoromethyl)phenoxy]-1-butenyl]cyclopentyl]-, [1 α (Z), 2 β (1E,3R*), 3 α , 5 α]-(+)-

OTHER NAMES:

CN (+)-Fluprostenol

CN 5-Heptenoic acid, 7-[3,5-dihydroxy-2-[3-hydroxy-4-[3-(trifluoromethyl)phenoxy]-1-butenyl]cyclopentyl]-, [1 α (Z), 2 β (1E,3R*), 3 α , 5 α]-

CN Fluoprostenol

CN Fluprostenol

FS STEREOSEARCH

DR 53468-75-0

MF C23 H29 F3 O6

CI COM

LC STN Files: AGRICOLA, ANABSTR, BEILSTEIN*, BIOBUSINESS, BIOSIS, BIOTECHNO, CA, CANCERLIT, CAPLUS, CHEMCATS, CHEMLIST, CSCHEM, DDFU, DRUGU, EMBASE, IFICDB, IFIPAT, IFIUDB, IPA, MEDLINE, MRCK*, TOXCENTER, USAN, USPAT2, USPATFULL, VETU

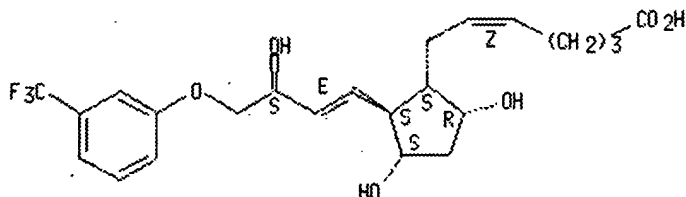
(*File contains numerically searchable property data)

Other Sources: EINECS**, WHO

(**Enter CHEMLIST File for up-to-date regulatory information)

Relative stereochemistry.

Double bond geometry as shown.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

116 REFERENCES IN FILE CA (1907 TO DATE)
 5 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA
 116 REFERENCES IN FILE CAPLUS (1907 TO DATE)

=> s chloprostenol

L2 0 CHLOPROSTENOL

=> s cloprostenol

L3 20 CLOPROSTENOL

=> d 20

L3 ANSWER 20 OF 20 REGISTRY COPYRIGHT 2003 ACS on STN

RN 40665-92-7 REGISTRY

CN 5-Heptenoic acid, 7-[(1R,2R,3R,5S)-2-[(1E,3R)-4-(3-chlorophenoxy)-3-hydroxy-1-butenyl]-3,5-dihydroxycyclopentyl]-, (5Z)-rel- (9CI) (CA INDEX NAME)

OTHER CA INDEX NAMES:

CN 5-Heptenoic acid, 7-[2-[4-(3-chlorophenoxy)-3-hydroxy-1-butenyl]-3,5-dihydroxycyclopentyl]-, [1α(Z),2β(1E,3R*),3α,5α]- (±)-

OTHER NAMES:

CN (±)-Cloprostenol

CN 5-Heptenoic acid, 7-[2-[4-(3-chlorophenoxy)-3-hydroxy-1-butenyl]-3,5-dihydroxycyclopentyl]-, [1α(Z),2β(1E,3R*),3α,5α]-

CN Cloprostenol

CN Estrofan

CN Estrophan

CN Estrophane

CN Oestrophan

CN Oestrophane

CN Racemic cloprostenol

FS STEREOSEARCH

DR 53529-41-2, 87347-50-0, 100786-10-5

MF C22 H29 Cl O6

CI COM

LC STN Files: AGRICOLA, ANABSTR, BEILSTEIN*, BIOBUSINESS, BIOSIS, BIOTECHNO, CA, CABA, CANCERLIT, CAPLUS, CASREACT, CBNB, CHEMCATS, CHEMLIST, CSCHM, DDFU, DRUGU, EMBASE, IFICDB, IFIPAT, IFIUDB, IPA, MEDLINE, MRCK*, PROMT, RTECS*, TOXCENTER, USAN, USPATFULL, VETU
 (*File contains numerically searchable property data)

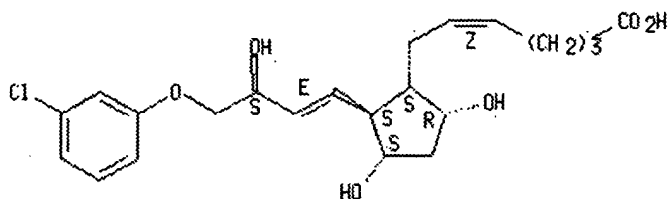
Other Sources: EINECS**, WHO

(**Enter CHEMLIST File for up-to-date regulatory information)

Relative stereochemistry.

STN Columbus

Double bond geometry as shown.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

610 REFERENCES IN FILE CA (1907 TO DATE)

6 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA

610 REFERENCES IN FILE CAPLUS (1907 TO DATE)

=>